

10/533,377

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NEWS 3 JUL 02 SCISEARCH enhanced with complete author names
NEWS 4 JUL 02 CHEMCATS accession numbers revised
NEWS 5 JUL 02 CA/CAplus enhanced with utility model patents from China
NEWS 6 JUL 16 CAplus enhanced with French and German abstracts
NEWS 7 JUL 18 CA/CAplus patent coverage enhanced
NEWS 8 JUL 26 USPATFULL/USPAT2 enhanced with IPC reclassification
NEWS 9 JUL 30 USGENE now available on STN
NEWS 10 AUG 06 CAS REGISTRY enhanced with new experimental property tags
NEWS 11 AUG 06 FSTA enhanced with new thesaurus edition
NEWS 12 AUG 13 CA/CAplus enhanced with additional kind codes for granted patents
NEWS 13 AUG 20 CA/CAplus enhanced with CAS indexing in pre-1907 records
NEWS 14 AUG 27 Full-text patent databases enhanced with predefined patent family display formats from INPADOCDB
NEWS 15 AUG 27 USPATOLD now available on STN
NEWS 16 AUG 28 CAS REGISTRY enhanced with additional experimental spectral property data
NEWS 17 SEP 07 STN AnaVist, Version 2.0, now available with Derwent World Patents Index
NEWS 18 SEP 13 FORIS renamed to SOFIS
NEWS 19 SEP 13 INPADOCDB enhanced with monthly SDI frequency
NEWS 20 SEP 17 CA/CAplus enhanced with printed CA page images from 1967-1998
NEWS 21 SEP 17 CAplus coverage extended to include traditional medicine patents
NEWS 22 SEP 24 EMBASE, EMBAL, and LEMBASE reloaded with enhancements
NEWS 23 OCT 02 CA/CAplus enhanced with pre-1907 records from Chemisches Zentralblatt
NEWS 24 OCT 19 BEILSTEIN updated with new compounds

NEWS EXPRESS 19 SEPTEMBER 2007: CURRENT WINDOWS VERSION IS V8.2, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 19 SEPTEMBER 2007.

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STRUCTURE FILE UPDATES: 9 NOV 2007 HIGHEST RN 952797-35-2
DICTIONARY FILE UPDATES: 9 NOV 2007 HIGHEST RN 952797-35-2

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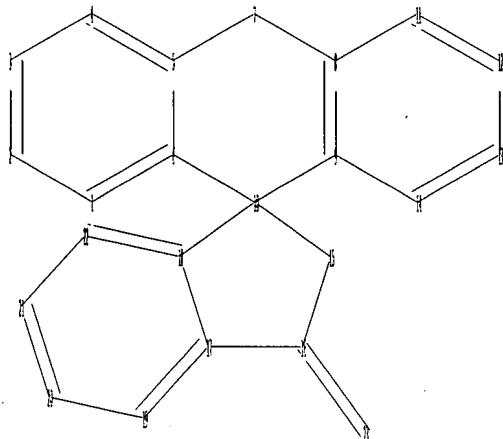
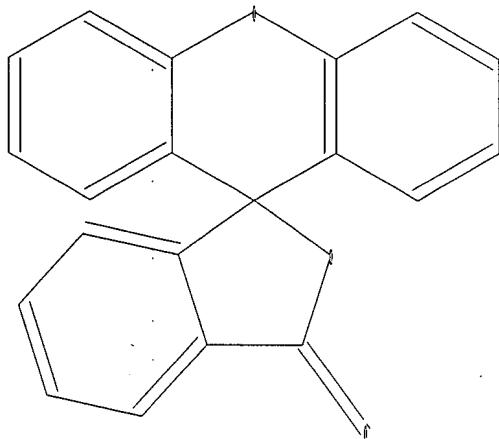
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<http://www.cas.org/support/stngen/stndoc/properties.html>

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=> Uploading C:\Program Files\Stnexp\Queries\10533377.str
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10/533,377



chain nodes :

23

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21

22

chain bonds :

16-23

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-10 7-8 8-9 8-11 9-10 9-14 10-15
10-18 11-12 12-13 13-14 15-16 16-17 17-18 17-19 18-22 19-20 20-21

21-22

exact/norm bonds :

16-23

exact bonds :

5-7 6-10 7-8 9-10 10-15 10-18 15-16 16-17

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6 8-9 8-11 9-14 11-12 12-13 13-14 17-18
17-19 18-22 19-20 20-21 21-22

isolated ring systems :

containing 1 :

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom
10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom
18:Atom 19:Atom 20:Atom 21:Atom 22:Atom 23:CLASS

L1 STRUCTURE UPLOADED

=> s 11

SAMPLE SEARCH INITIATED 21:21:34 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 833 TO ITERATE

10/533,377

100.0% PROCESSED 833 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

50 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 14929 TO 18391
PROJECTED ANSWERS: 13282 TO 16558

L2 50 SEA SSS SAM L1

=> s 11 ful
FULL SEARCH INITIATED 21:21:41 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 16121 TO ITERATE

100.0% PROCESSED 16121 ITERATIONS
SEARCH TIME: 00.00.01

14354 ANSWERS

L3 14354 SEA SSS FUL L1 |

=> file caplus
COST IN U.S. DOLLARS | SINCE FILE | TOTAL
FULL ESTIMATED COST | ENTRY | SESSION
172.10 | 172.31

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FILE LAST UPDATED: 9 Nov 2007 (20071109/ED)

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=> s 13
L4 33637 L3

10/533,377

=> s 14 and (prepar? or method or make or making or synthe? or process)
1828766 PREPAR?
133837 PREP
2315 PREPS
135936 PREP
(PREP OR PREPS)
2125686 PREPD
3 PREPDS
2125688 PREPD
(PREPD OR PREPDS)
146675 PREPG
9 PREPGS
146683 PREPG
(PREPG OR PREPGS)
2858263 PREPN
211266 PREPNS
3017389 PREPN
(PREPN OR PREPNS)
5074260 PREPAR?
(PREPAR? OR PREP OR PREPD OR PREPG OR PREPN)
3548243 METHOD
1418435 METHODS
4566798 METHOD
(METHOD OR METHODS)
273407 MAKE
212582 MAKES
470609 MAKE
(MAKE OR MAKES)
329221 MAKING
35 MAKINGS
329250 MAKING
(MAKING OR MAKINGS)
1670374 SYNTHE?
2518307 PROCESS
1716032 PROCESSES
3755651 PROCESS
(PROCESS OR PROCESSES)
L5 18833 L4 AND (PREPAR? OR METHOD OR MAKE OR MAKING OR SYNTHE? OR
PROCE
SS)
=> s 15 and esther
71 ESTHER
11 ESTHERS
82 ESTHER
(ESTHER OR ESTHERS)
L6 0 L5 AND ESTHER
=> s 15 and ester
610455 ESTER
448187 ESTERS

10/533,377

846715 ESTER
(ESTER OR ESTERS)

L7 1496 L5 AND ESTER

=> s 15 and acid ester

4474144 ACID
1600200 ACIDS
4978785 ACID
(ACID OR ACIDS)
610455 ESTER
448187 ESTERS
846715 ESTER
(ESTER OR ESTERS)

149136 ACID ESTER
(ACID(W)ESTER)

L8 208 L5 AND ACID ESTER

=> s 18 and (phthalein or phthalacein or fluorescein)

877 PHTHALEIN
357 PHTHALEINS
1091 PHTHALEIN
(PHTHALEIN OR PHTHALEINS)
0 PHTHALACEIN
27568 FLUORESCEIN
259 FLUORESCEINS
27645 FLUORESCEIN
(FLUORESCEIN OR FLUORESCEINS)

L9 49 L8 AND (PHTHALEIN OR PHTHALACEIN OR FLUORESCEIN)

L

L9 ANSWER 24 OF 49 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 1992:241720 CAPLUS

DOCUMENT NUMBER: 116:241720

TITLE: Product for improved permanent waving of hair and simultaneously coloring and permanently waving hair

INVENTOR(S): Schultz, Thomas M.; Patel, Jitendra; Wong,

Stephanie

PATENT ASSIGNEE(S): Shiseido Co., Ltd., Japan

SOURCE: U.S., 6 pp.

CODEN: USXXAM

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 5094662	A	19920310	US 1990-612227	19901113
US 5188639	A	19930223	US 1992-847788	19920306
PRIORITY APPLN. INFO.:			US 1990-612227	A3 19901113

AB By combining of a fluorescein-based dye with a mercaptan-based permanent waving composition and maintaining the pH of the resulting composition between .apprx.2.5-4.5, a composition is achieved which simultaneously colors and permanently waves the hair. In addition, even if no coloring or dyeing of the hair is desired, the use of a colorless or complimentary fluorescein-based dye with the mercaptan-based permanent waving composition achieves a composition which imparts a substantially improved curl configuration to the hair as well as substantially longer lasting curls. Preferably, the mercaptan-based permanent waving composition employed comprises an ester of either thioglycolic acid, thiolactic acid, or the amide of 2-aminoethanethiol. Hair treated with 1 part 32.5% glycerin monothioglycolate in glycerin; 1 part 1.50% NH4Cl in deionized water, 0.005 part D & C Red Number 28 5.0% in deionized water, and preservative q.s.; pH 3.2, had improved curl retention and structural integrity.

Hair

color was dark auburn.

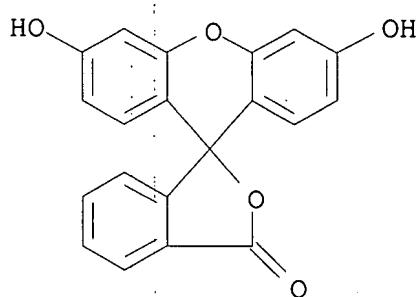
IT 2321-07-5D, Fluorescein, derivs. 15086-94-9
17372-87-1, D&C Red Number 22 18472-87-2

RL: BIO (Biological study)

(permanent wave and hair dye preparation containing)

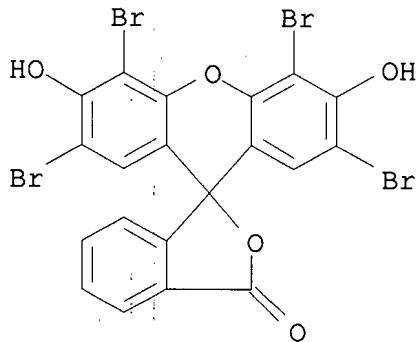
RN 2321-07-5 CAPLUS

CN Spiro[isobenzofuran-1(3H),9'-[9H]xanthen]-3-one, 3',6'-dihydroxy- (CA INDEX NAME)



RN 15086-94-9 CAPLUS

CN Spiro[isobenzofuran-1(3H),9'-[9H]xanthen]-3-one,
2',4',5',7'-tetrabromo-
3',6'-dihydroxy- (CA INDEX NAME)

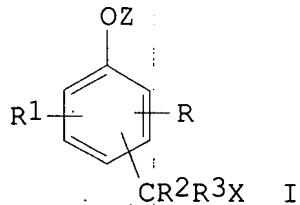


RN 17372-87-1 CAPLUS
 CN Spiro[isobenzofuran-1(3H),9'-[9H]xanthen]-3-one,
 2',4',5',7'-tetrabromo-
 3',6'-dihydroxy-, sodium salt (1:2) (CA INDEX NAME)

L9 ANSWER 27 OF 49 CAPLUS COPYRIGHT 2007 ACS on STN
 ACCESSION NUMBER: 1990:511969 CAPLUS
 DOCUMENT NUMBER: 113:111969
 TITLE: Enzyme-controlled-release system using a
 use quinone-methide elimination reaction mechanism for
 in immunoassays and pharmaceuticals
 INVENTOR(S): Meneghini, Frank A.; Palumbo, Paul S.
 PATENT ASSIGNEE(S): Polaroid Corp., USA
 SOURCE: PCT Int. Appl., 34 pp.
 CODEN: PIXXD2
 DOCUMENT TYPE: Patent
 LANGUAGE: English
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 9001558	A1	19900222	WO 1989-US1696	19890420
W: JP				
RW: DE, FR, GB, IT, NL				
US 5112739	A	19920512	US 1988-227141	19880802
EP 396642	A1	19901114	EP 1989-907960	19890420
EP 396642	B1	19940511		
R: DE, FR, GB, IT, NL				
JP 03500367	T	19910131	JP 1989-507309	19890420
CA 1336586	C	19950808	CA 1989-598563	19890503
PRIORITY APPLN. INFO.:			US 1988-227141	A 19880802
			WO 1989-US1696	W 19890420

OTHER SOURCE(S): MARPAT 113:111969
GI



AB An enzyme-controlled-release system uses compound I (R, R1, R2, R3 = H, substituent affecting the mobility or reactivity of the compound, or a substituent including a biol. active group; X = leaving group and may be

an organic, organometallic, or inorg. moiety; Z = enzyme substrate cleavable

by an active enzyme; CR2R3X is either ortho or para to the OZ moiety).

An

active enzyme cleaves the substrate, Z; the resultant active intermediate

undergoes a quinone-methide elimination reaction to release the leaving group X. The system is useful for detecting an analyte of interest and may be used in, e.g., immunoassays, enzyme amplification systems, and the

release of pharmacol. active ligands.

(4-Resorufinylmethyl-2-nitrophenyl)-

2,3,4,6-tetra-O-acetyl- β -D-galactopyranoside was prepared by heating a solution of

(4-chloromethyl-2-nitrophenyl)-2,3,4,6-tetra-O-acetyl-

β -D-galactopyranoside (preparation given), Na resorufin, and a catalytic amount of NaI in dry DMF at 70° for 4 h. The galactosyl acetate protecting groups were removed with NaOMe. When the galactopyranoside was treated with β -galactosidase, the leaving group release rate was 0.25 (compared with 1.0 for o-nitrophenolgalactoside).

IT 2321-07-5, Fluorescein 2321-07-5D,

Fluorescein, derivs. 3086-44-0D, Rhodol, derivs.

RL: ANST (Analytical study)

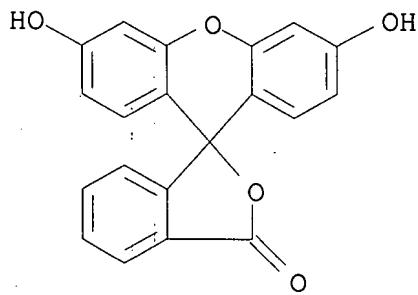
(enzyme-controlled-release compound containing, quinone-methide elimination

in, for immunochem. anal.)

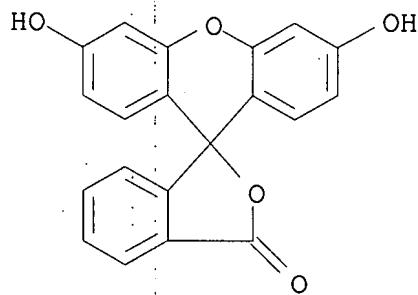
RN 2321-07-5 CAPLUS

CN Spiro[isobenzofuran-1(3H),9'-[9H]xanthen]-3-one, 3',6'-dihydroxy- (CA INDEX NAME)

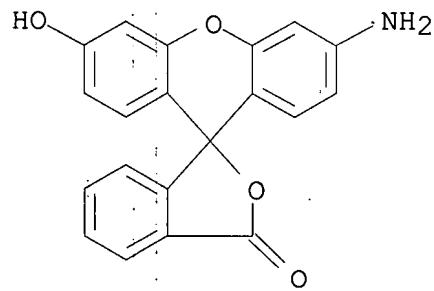
10/533,377



RN 2321-07-5 CAPLUS
CN Spiro[isobenzofuran-1(3H),9'-[9H]xanthen]-3-one, 3',6'-dihydroxy- (CA
INDEX NAME)



RN 3086-44-0 CAPLUS
CN Spiro[isobenzofuran-1(3H),9'-[9H]xanthen]-3-one, 3'-amino-6'-hydroxy-
(CA
INDEX NAME)



L9 ANSWER 28 OF 49 CAPLUS COPYRIGHT 2007 ACS on STN
ACCESSION NUMBER: 1989:570653 CAPLUS
DOCUMENT NUMBER: 111:170653
TITLE: Novel amphiphilic nucleic acid conjugates with more
efficient membrane transport, their
preparation and use

INVENTOR(S): Tullis, Richard H.
 PATENT ASSIGNEE(S): Synthetic Genetics, USA
 SOURCE: PCT Int. Appl., 48 pp.
 CODEN: PIXXD2
 DOCUMENT TYPE: Patent
 LANGUAGE: English
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

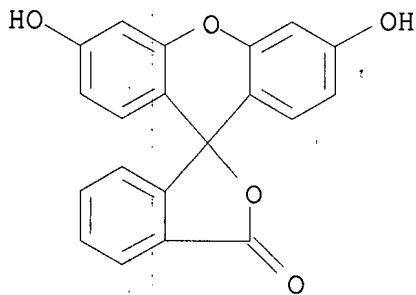
PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 8809810	A1	19881215	WO 1988-US2009	19880611
W: JP				
RW: AT, BE, CH, DE, FR, GB, IT, LU, NL, SE				
US 4904582	A	19900227	US 1987-61874	19870611
EP 321548	A1	19890628	EP 1988-906384	19880611
R: AT, BE, CH, DE, FR, GB, IT, LI, LU, NL, SE				
JP 03500530	T	19910207	JP 1988-505633	19880611
PRIORITY APPLN. INFO.:			US 1987-61874	A 19870611
			WO 1988-US2009	W 19880611

AB Novel nucleic acid conjugates are prepared comprising a relatively short nucleic acid sequence complementary to a sequence of interest for modifying intracellular expression, a linking group, and a group imparting amphiphilic character, usually more hydrophobic than hydrophilic. The resulting conjugates are more efficient in membrane transport and can cross the membrane and effectively modulate the transcriptional system. Oligonucleotides 20 bases long (synthesized antisense to mouse β -globin mRNA) were conjugated at the 5'-terminal to PEG using imidazole-activated carboxylic acid esters and bis-aminoalkyl PEG. At 15 μ M the conjugate selectively inhibited Hb synthesis in cultured Friend murine erythroleukemia cells by 95% compared to 0%, 24%, and 78% inhibition with DMSO, PEG, and PEG + the oligonucleotide, resp., at 100 μ M.

IT 2321-07-5D, Fluorescein, oligonucleotide conjugates
 RL: ANST (Analytical study)
 (mRNA maturation or translation inhibition by)

RN 2321-07-5 CAPLUS

CN Spiro[isobenzofuran-1(3H),9'-[9H]xanthen]-3-one, 3',6'-dihydroxy- (CA INDEX NAME)

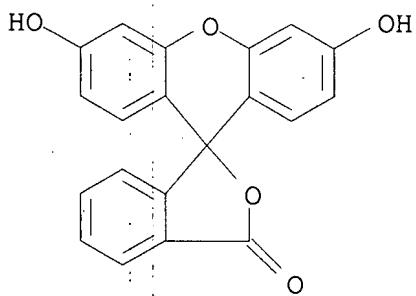


IT 27072-45-3, FITC

RL: RCT (Reactant); RACT (Reactant or reagent)
(reaction of, with diamine and oligonucleotide)

RN 27072-45-3 CAPLUS

CN Spiro[isobenzofuran-1(3H),9'-[9H]xanthen]-3-one, 3',6'-dihydroxy-5(or 6)-isothiocyanato- (CA INDEX NAME)



D1-N=C=S

L9 ANSWER 29 OF 49 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 1988:551435 CAPLUS

DOCUMENT NUMBER: 109:151435

TITLE: The characterization of metal-containing polymeric dyes for control of polymer degradation

AUTHOR(S): Carragher, Charles E., Jr.; Linville, Raymond J.; Stevenson, Donald F.; Foster, Van R.; Williams, Melanie; Alois, Mary Jo

CORPORATE SOURCE: Dep. Chem., Florida Atlantic Univ., Boca Raton, FL, 33431, USA

SOURCE: Polymeric Materials Science and Engineering (1988), 58, 85-9

CODEN: PMSEDG; ISSN: 0743-0515

DOCUMENT TYPE: Journal

LANGUAGE: English

AB The title dyes were prepared by copolymerg. R2SnCl2 (R = Me, Et, Bu,

Ph, cyclohexyl) with eosin Y, mercurochrome, phloxine B, eosin B, fluorescein, or Rose Bengal and tested for inhibition of *A. niger* in acrylic latex pastes. Polymeric dyes prepared from biscyclopentadienyltitanium dichloride were also discussed.

IT 89761-57-9 116828-87-6 116828-90-1
116828-91-2 116852-57-4 116852-58-5

RL: USES (Uses)

(bactericidal properties of colored)

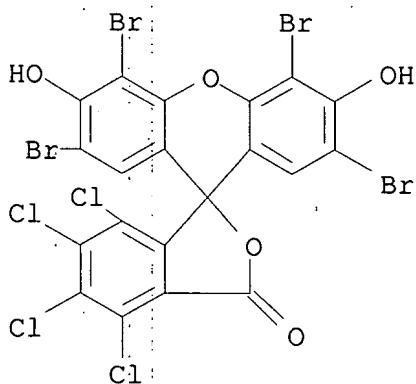
RN 89761-57-9 CAPLUS

CN Spiro[isobenzofuran-1(3H),9'-[9H]xanthen]-3-one,
2',4',5',7'-tetrabromo-
4,5,6,7-tetrachloro-3',6'-dihydroxy-, disodium salt, polymer with
dichlorodiphenylstannane (9CI) (CA INDEX NAME)

CM 1

CRN 18472-87-2

CMF C20 H4 Br4 Cl14 O5 . 2 Na

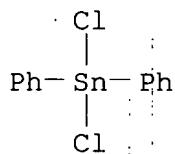


●2 Na

CM 2

CRN 1135-99-5

CMF C12 H10 Cl12 Sn



10/533,377

RN 116828-87-6 CAPLUS

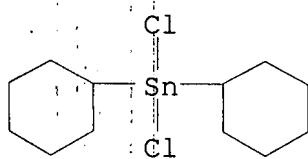
CN Mercury,

(2',7'-dibromo-3',6'-dihydroxy-3-oxospiro[isobenzofuran-1(3H),9'-[9H]xanthen]-4'-yl)hydroxy-, disodium salt, polymer with dichlorodicyclohexylstannane (9CI) (CA INDEX NAME)

CM 1

CRN 3342-69-6

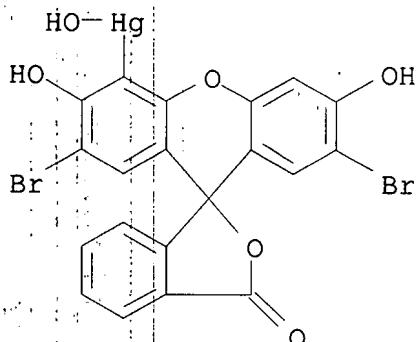
CMF C12 H22 Cl2 Sn



CM 2

CRN 129-16-8

CMF C20 H10 Br2 Hg O6 . 2 Na



●2 Na

RN 116828-90-1 CAPLUS

CN Mercury,

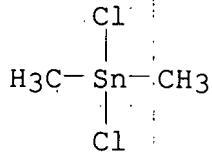
(2',7'-dibromo-3',6'-dihydroxy-3-oxospiro[isobenzofuran-1(3H),9'-[9H]xanthen]-4'-yl)hydroxy-, disodium salt, polymer with dichlorodimethylstannane (9CI) (CA INDEX NAME)

CM 1

CRN 753-73-1

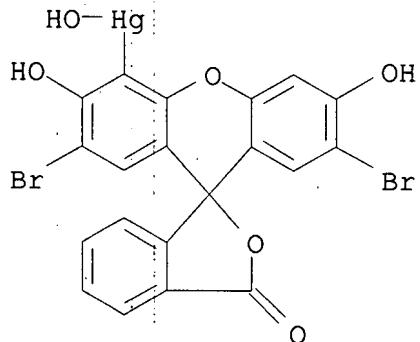
10/533,377

CMF C2 H6 Cl2 Sn



CM 2

CRN 129-16-8
CMF C20 H10 Br2 Hg O6 . 2 Na

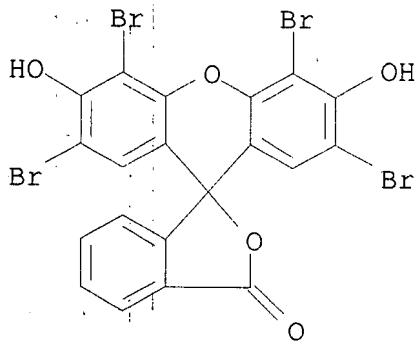


●2 Na

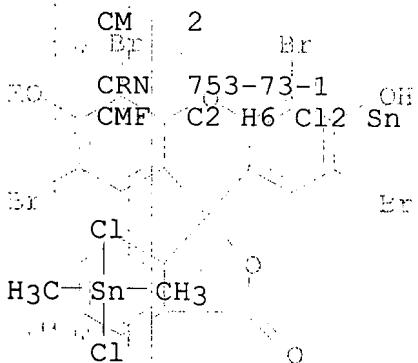
RN 116828-91-2 CAPLUS
CN Spiro[isobenzofuran-1(3H),9'-[9H]xanthen]-3-one,
2',4',5',7'-tetrabromo-
3',6'-dihydroxy-, disodium salt, polymer with dichlorodimethylstannane
(9CI) (CA INDEX NAME)

CM 1

CRN 17372-87-1
CMF C20 H8 Br4 O5 . 2 Na



10/533,377-2 Na



RN 116852-57-4 CAPLUS

CN Mercury,

(2',7'-dibromo-3',6'-dihydroxy-3-oxospiro[isobenzofuran-1(3H),9'-[9H]xanthen]-4'-yl)hydroxy-, disodium salt, polymer with dibutyltin dichloride (9CI) (CA INDEX NAME)

CM 1

CBM 116852-57-4

CRN 683-18-1

CMF C8 H18 Cl2 Sn

R:

Cl

n-Bu-Sn-Bu-n

Cl

RN 116852-57-4 CAPLUS

(2',7'-dibromo-3',6'-dihydroxy-3-oxospiro[isobenzofuran-1(3H),9'-[9H]xanthen]-4'-yl)hydroxy-, disodium salt, polymer with dibutyltin dichloride (9CI) (CA INDEX NAME)

CM 1

CBM 116852-57-4

CRN 683-18-1

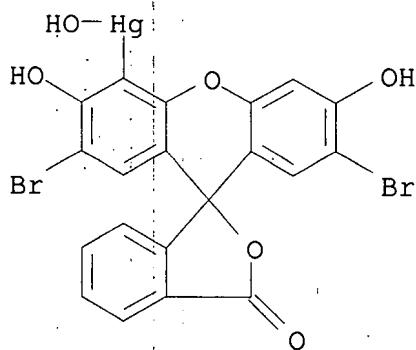
CMF C8 H18 Cl2 Sn

10/533/377

CM 2

CRN 129-16-8

CMF C20 H10 Br2 Hg O6 . 2 Na



●2 Na

RN 116852-58-5 CAPLUS

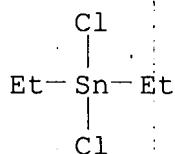
CN Mercury,

(2',7'-dibromo-3',6'-dihydroxy-3-oxospiro[isobenzofuran-1(3H),9'-[9H]xanthen]-4'-yl)hydroxy-, disodium salt, polymer with dichlorodiethylstannane (9CI) (CA INDEX NAME)

CM 1

CRN 866-55-7

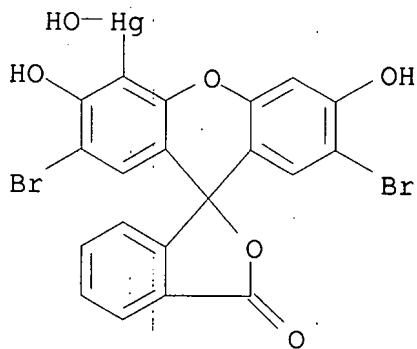
CMF C4 H10 Cl2 Sn



CM 2

CRN 129-16-8

CMF C20 H10 Br2 Hg O6 . 2 Na



●2 Na

IT 89761-58-0 89777-76-4 116828-89-8

RL: USES (Uses)

(testing of colored, for bactericidal properties)

RN 89761-58-0 CAPLUS

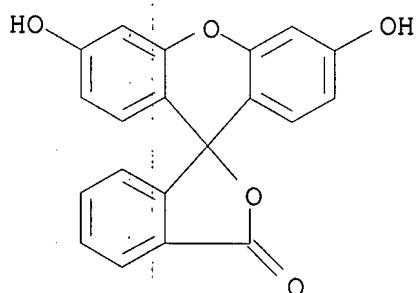
CN Spiro[isobenzofuran-1(3H),9'-[9H]xanthen]-3-one, 3',6'-dihydroxy-, polymer

with dichlorodiphenylstannane (9CI) (CA INDEX NAME)

CM 1

CRN 2321-07-5

CMF C20 H12 O5

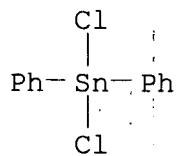


CM 2

CRN 1135-99-5

CMF C12 H10 Cl2 Sn

10/533,377



RN 89777-76-4 CAPLUS

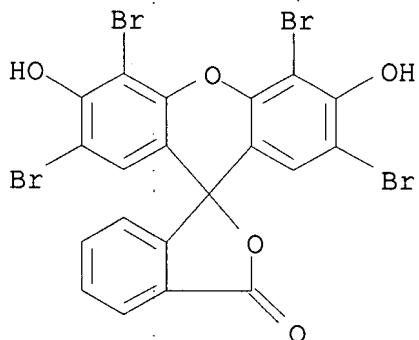
CN Spiro[isobenzofuran-1(3H),9'-[9H]xanthen]-3-one,
2',4',5',7'-tetrabromo-

3',6'-dihydroxy-, disodium salt, polymer with dichlorodiphenylstannane
(9CI) (CA INDEX NAME)

CM 1

CRN 17372-87-1

CMF C20 H8 Br4 O5 . 2 Na

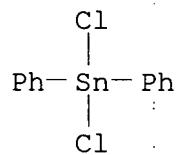


●2 Na

CM 2

CRN 1135-99-5

CMF C12 H10 Cl2 Sn



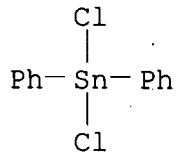
RN 116828-89-8 CAPLUS

10/533,377

CN Spiro[isobenzofuran-1(3H),9'-[9H]xanthen]-3-one, 4',5'-dibromo-3',6'-dihydroxy-2',7'-dinitro-, disodium salt, polymer with dichlorodiphenylstannane (9CI) (CA INDEX NAME)

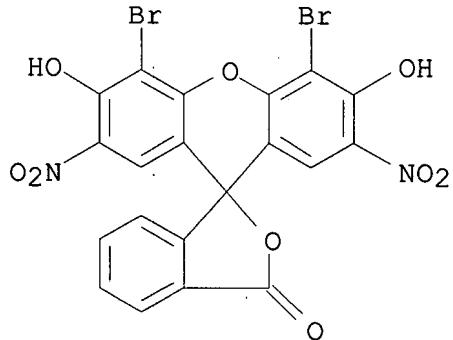
CM 1

CRN 1135-99-5
CMF C12 H10 Cl2 Sn



CM 2

CRN 548-24-3
CMF C20 H8 Br2 N2 O9 . 2 Na



●2 Na

L9 ANSWER 33 OF 49 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 1972:444650 CAPLUS

DOCUMENT NUMBER: 77:44650

ORIGINAL REFERENCE NO.: 77:7391a,7394a

TITLE: Hydrolysis of phthalyl amino acid
esters of fluorescein in the
presence of leucine aminopeptidase

AUTHOR(S): Thomas, John J.; Eveland, Warren C.; Medzon, Edward
L.; Christian, Walter; Wylie, Dwayne E.;

Burckhalter,

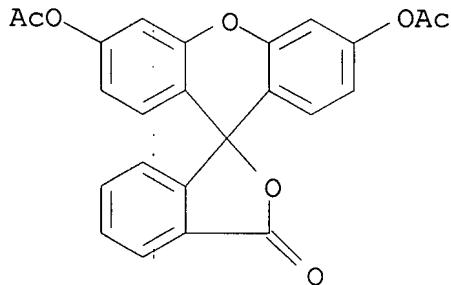
CORPORATE SOURCE: Joseph; Jones, Ronald H.
 SOURCE: Dep. Med. Chem., Univ. Michigan, Ann Arbor, MI, USA
 Proceedings of the Society for Experimental Biology
 and Medicine (1972), 140(1), 179-82
 CODEN: PSEBAA; ISSN: 0037-9727

DOCUMENT TYPE: Journal
 LANGUAGE: English

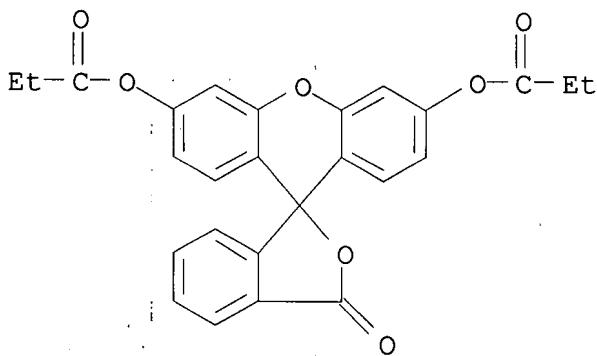
GI For diagram(s), see printed CA Issue.
 AB Phthalyl amino acid esters of fluorescein were synthesized and characterized. They were tested (in vitro) for use as possible fluorogenic substrates for leucine aminopeptidase (LAP). The rates of enzymic hydrolysis of these compds. were compared with those of some com. available fluorescein esters. The phthalyl amino acid esters were hydrolyzed rapidly by LAP, lipase, and chymotrypsin and much more slowly by trypsin and cholinesterases. The min. detectable LAP concentration was 1 μ g/ml for d,
 d-bis(α -isopropyl-1,3-dioxo- 2-isoindolinylacetyl) fluorescein (d,d-I) to 8 μ g/ml for the 1,1-bis(α -isobutyl- 1,3-dioxo- 2-isoindolinylacetyl)-(1,1-II) and 1,1-bis(α -butyl-1,3-dioxo-2-isoindolinylacetyl) fluorescein (1,1-III).

IT 596-09-8 7276-28-0 7298-65-9 7308-90-9
 19722-86-2 36889-44-8 36889-45-9
 36889-46-0 36889-47-1 36889-48-2
 36905-09-6 36984-36-8
 RL: BIOL (Biological study)
 (hydrolysis by leucine aminopeptidase)

RN 596-09-8 CAPLUS
 CN Spiro[isobenzofuran-1(3H),9'-[9H]xanthen]-3-one, 3',6'-bis(acetyloxy)-
 (CA INDEX NAME)



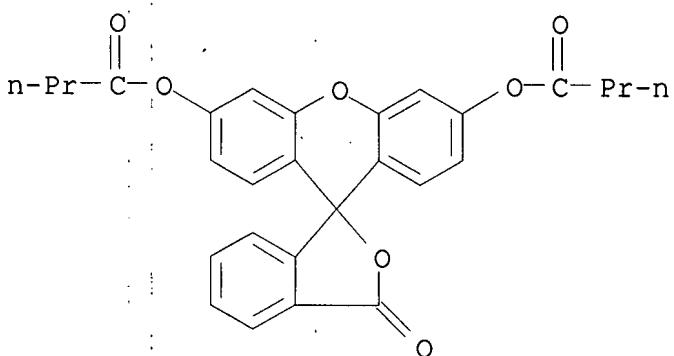
RN 7276-28-0 CAPLUS
 CN Spiro[isobenzofuran-1(3H),9'-[9H]xanthen]-3-one,
 3',6'-bis(1-oxoproxy)-
 (CA INDEX NAME)



RN 7298-65-9 CAPLUS

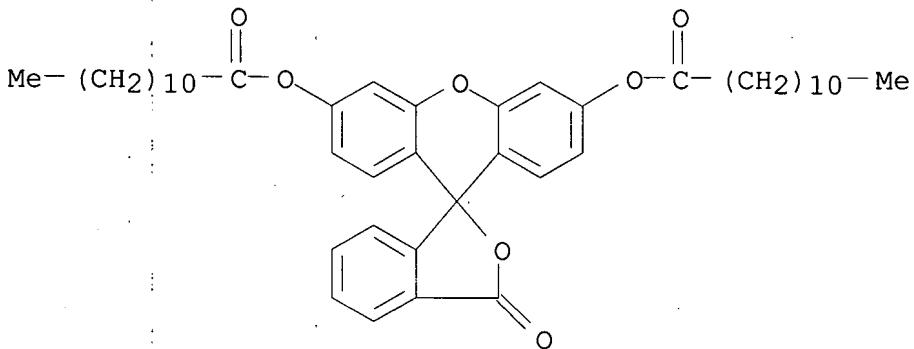
CN Butanoic acid,

1,1'-(3-oxospiro[isobenzofuran-1(3H),9'-[9H]xanthene]-3',6'-diyl) ester (CA INDEX NAME)



RN 7308-90-9 CAPLUS

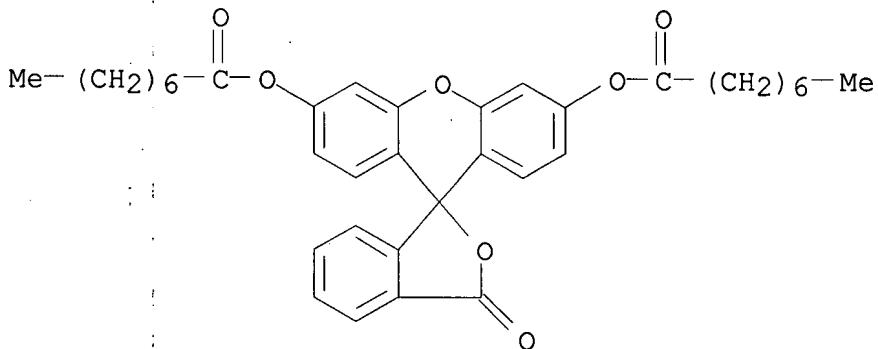
CN Dodecanoic acid, 1,1'-(3-oxospiro[isobenzofuran-1(3H),9'-[9H]xanthene]-3',6'-diyl) ester (CA INDEX NAME)



RN 19722-86-2 CAPLUS

10/533,377

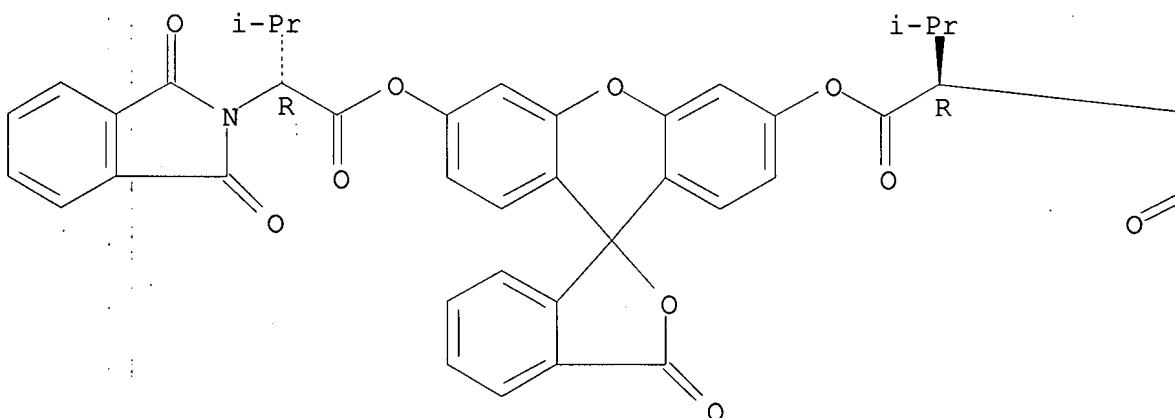
CN Octanoic acid,
1,1'-[3-oxospiro[isobenzofuran-1(3H),9'-[9H]xanthene]-3',6'-diyl] ester (CA INDEX NAME)



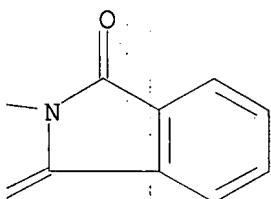
RN 36889-44-8 CAPLUS
CN 2H-Isoindole-2-acetic acid, 1,3-dihydro- α -(1-methylethyl)-1,3-dioxo-, 3-oxospiro[isobenzofuran-1(3H),9'-[9H]xanthene]-3',6'-diyl ester, (R*,R*)-(+)-(9CI) (CA INDEX NAME)

Rotation (+). Absolute stereochemistry unknown.

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PAGE 1-B



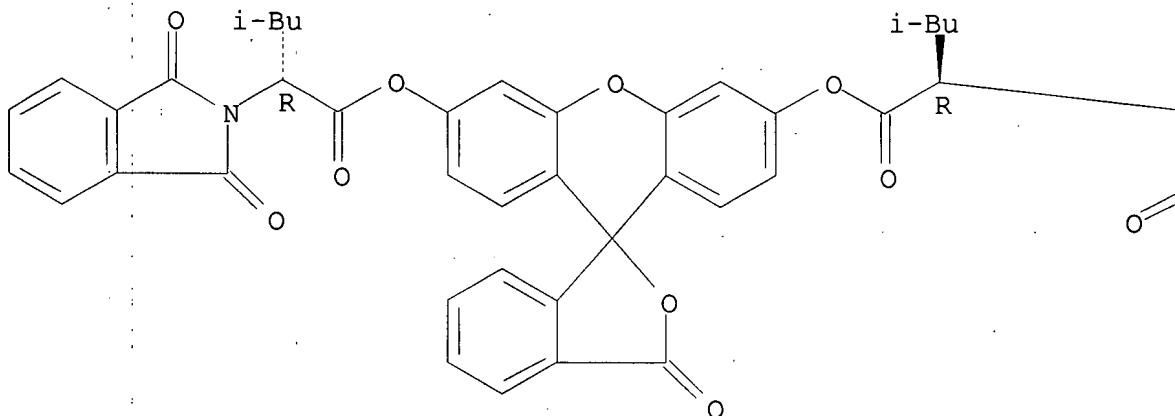
10/533,377

RN 36889-45-9 CAPLUS

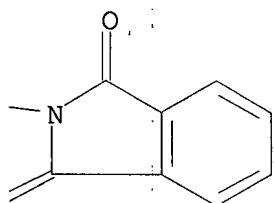
CN 2H-Isoindole-2-acetic acid, 1,3-dihydro- α -(2-methylpropyl)-1,3-dioxo-, 3-oxospiro[isobenzofuran-1(3H),9'-[9H]xanthene]-3',6'-diyl ester, (R*,R*)-(-)-(9CI) (CA INDEX NAME)

Rotation (-). Absolute stereochemistry unknown.

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PAGE 1-B

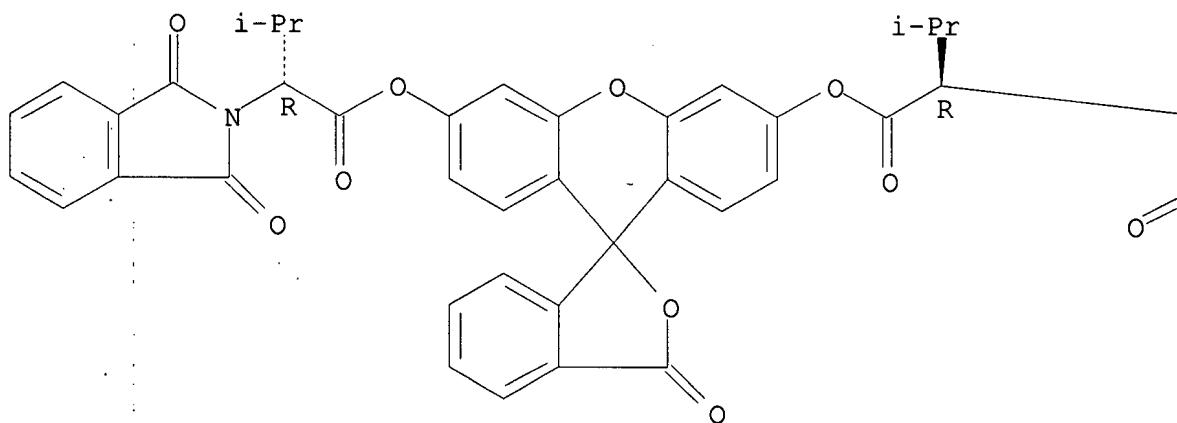


RN 36889-46-0 CAPLUS

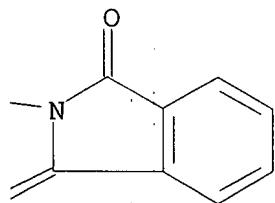
CN 2H-Isoindole-2-acetic acid, 1,3-dihydro- α -(1-methylethyl)-1,3-dioxo-, 3-oxospiro[isobenzofuran-1(3H),9'-[9H]xanthene]-3',6'-diyl ester, (R*,R*)-(-)-(9CI) (CA INDEX NAME)

Rotation (-). Absolute stereochemistry unknown.

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PAGE 1-B

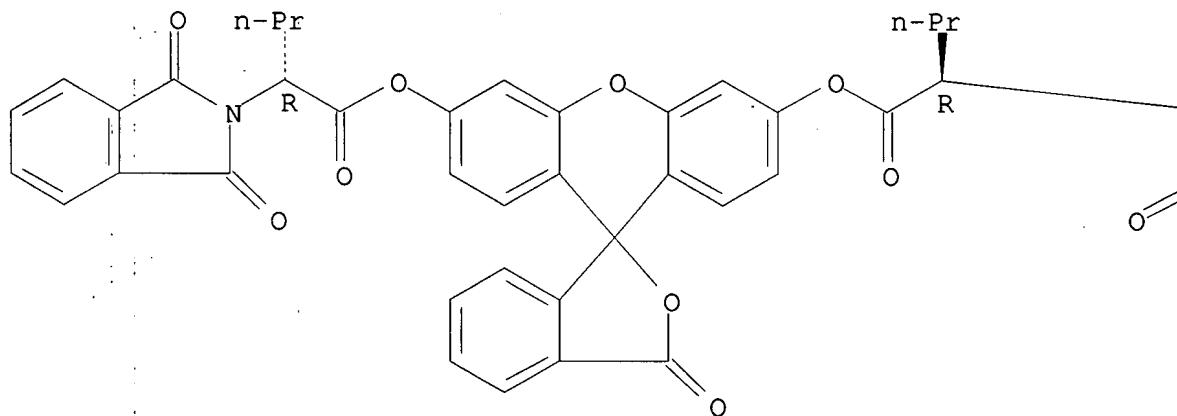


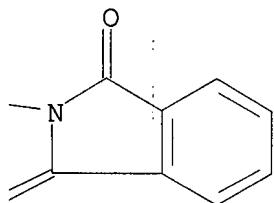
RN 36889-47-1 CAPLUS

CN 2H-Isoindole-2-acetic acid, 1,3-dihydro-1,3-dioxo- α -propyl-, 3-oxospiro[isobenzofuran-1(3H),9'-[9H]xanthene]-3',6'-diyl ester, (R*,R*)-(-)- (9CI) (CA INDEX NAME)

Rotation (-). Absolute stereochemistry unknown.

PAGE 1-A

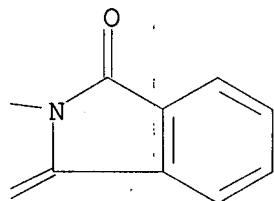
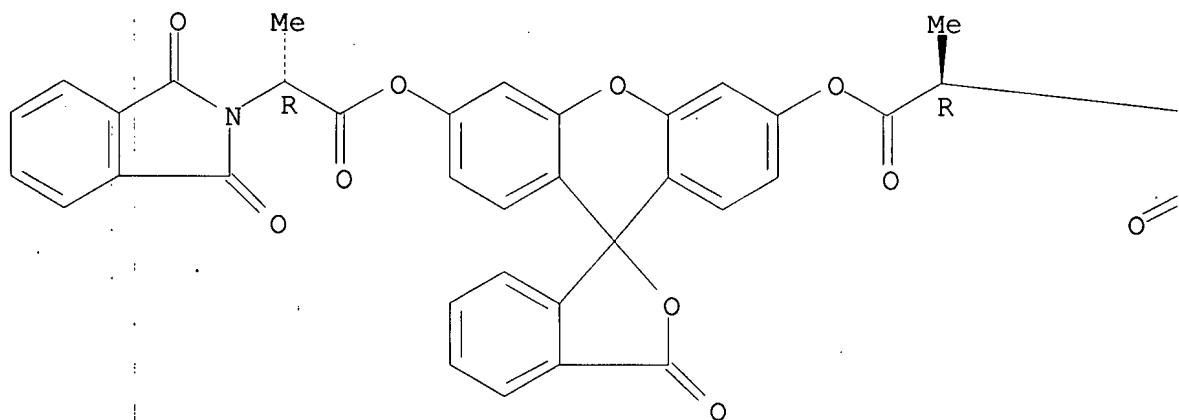




RN 36889-48-2 CAPLUS

CN 2H-Isoindole-2-acetic acid, 1,3-dihydro- α -methyl-1,3-dioxo-,
3-oxospiro[isobenzofuran-1(3H),9'-[9H]xanthene]-3',6'-diyl ester,
(R*,R*)-(-) - (9CI) (CA INDEX NAME)

Rotation (-). Absolute stereochemistry unknown.



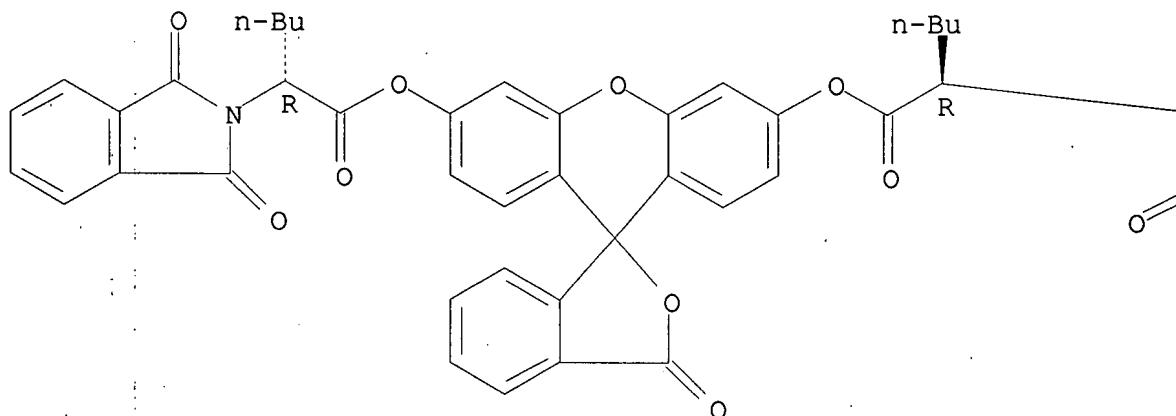
RN 36905-09-6 CAPLUS

CN 2H-Isoindole-2-acetic acid, α -butyl-1,3-dihydro-1,3-dioxo-,
3-oxospiro[isobenzofuran-1(3H),9'-[9H]xanthene]-3',6'-diyl ester,
(R*,R*)-(-) - (9CI) (CA INDEX NAME)

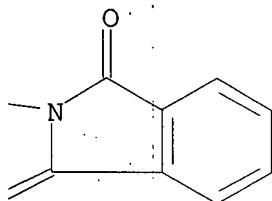
10/533,377

Rotation (-). Absolute stereochemistry unknown.

PAGE 1-A



PAGE 1-B

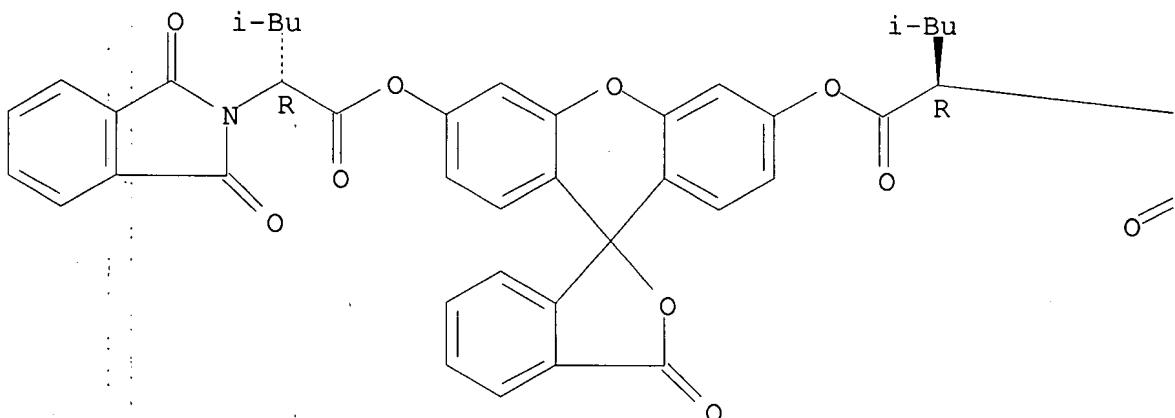


RN 36984-36-8 CAPLUS

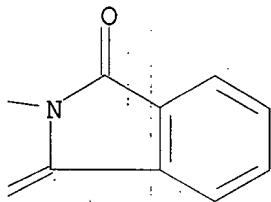
CN 2H-Isoindole-2-acetic acid, 1,3-dihydro- α -(2-methylpropyl)-1,3-dioxo-, 3-oxospiro[isobenzofuran-1(3H), 9'-[9H]xanthene]-3',6'-diyl ester, (R*,R*)-(+)- (9CI) (CA INDEX NAME)

Rotation (+). Absolute stereochemistry unknown.

PAGE 1-A



PAGE 1-B



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L9 ANSWER 49 OF 49 CAPLUS COPYRIGHT 2007 ACS on STN

ACCESSION NUMBER: 1925:16971 CAPLUS

DOCUMENT NUMBER: 19:16971

ORIGINAL REFERENCE NO.: 19:2196i,2197a-g

TITLE: The condensation of resorcinols and a few other aromatic hydroxy compounds with some acids, esters, lactones and lactams

AUTHOR(S): Sen, R. N.; Sircar, S. S.

SOURCE: Quart. J. Indian Chem. Soc. (1924), 1, 151-72

DOCUMENT TYPE: Journal

LANGUAGE: Unavailable

AB cf. C. A. 18, 1831. Hthalein-like compds. were obtained by heating 1 mol.

HO compound with 2 mols. acid or acid derivative and powdered ZnCl₂ 3 h. to180-200° in a current of dry HCl. Rhodamies were prepared at a lower temperature and without HCl. The product was freed from ZnCl₂ by

extraction with HCl, dissolved in alkali and precipitated with dilute HCl or AcOH.

Bromination was carried out by allowing the alc. solution to stand overnight

with a slight excess of Br. Esters react with phenols more readily than

the free acids. It is remarkable that pyrogallol derivs. are not fluorescent in alkaline and organic solns., unlike the resorcinol derivs. In

dibasic acids the 2 CO₂H groups can react successively. Simultaneous reaction can be forced only by drastic means and with poor yields.

Resorcinolgallein (80% yield), dark red, soluble in alkaline and organic solvents

with green-red fluorescence, insol. in AcOH, ether and C₆H₆, dyes brownish

shades on wool and silk, does not m. 250°. Di-Br derivative, red, non-fluorescent in alkali, dyes wool and silk deep red, does not m. 250°. Resorcinol-o-amino-benzein softens 1.75-7° and

closely resembles the salicylein in color, fluorescence, solubility and dyeing

properties. Yield 70%. Di-Br-derivative resembles the salicylein derivative,

decomps. 195° and dyes red shades. K salt dissolves in water with green-red fluorescence changing to bluish red on standing.

Resorcinolstearein softens 152°. The deep red color changes to a bluish tone on standing. The fluorescence is slight but appreciable, the

affinity for wool and silk slight. Resorcinolpyromucein, insol. in C₆H₆

and ether, similar to the stearein in color and fluorescence, has more pronounced dyeing properties, does not m. 250°.

Pyrogallolsalicylein, does not m. 250°; the deep red-brown alkaline and the organic solns. are non-fluorescent; it is soluble in acetone and a mixture of

C₆H₆N and water. Anthranilorhodamine, pink powder, soluble in acids with

green-red fluorescence, more marked in organic solns., dyes wool with a violet-greenish shade. and softens 230°. Yield 60%.

Resorcinolsalicylein, does not m. 260°, and pyrogallolsalicylein were prepared in 85% yield from Me salicylate.

β-Naphtholcoumarein a yellowish brown powder, soluble in NaOH, acetone,

AcOH and concentrated H₂SO₄ with green fluorescence, softens 115° and is

precipitated from alkaline solns. by CO₂. The solubility in NaOH is due to the opening of

the lactone ring, which is confirmed by anal. α-Naphtholcoumarein, a dark brown powder soluble in NaOH with red color and softening 117°, is obtained in 40% yield with considerable tarring. The slight fluorescence in NaOH is attributable to partial o-condensation to a fluoran like substance. Both compds. have a hardly appreciable affinity

for animal fibers. Pyrogallolcoumarein, analogous to other pyrogallol derivs., does not m. 250°. Coumarinrhodamine dissolves in acids

with violet color and greenish fluorescence more marked in organic solvents,

softens 156° and dyes wool and silk violet shades.

Phenolresorcinolphthalein, prepared in 70% yield from phenolphthalein and resorcinol by heating 4.5 h. to 210-5°, while 3-4 h. heating to 180-200° yields fluorescein, orange powder, somewhat soluble in hot water, resembles fluorescein in alkaline solution, does not m. 250° and dyes a greenish yellow shade.

The

tetra-Br-derivative is an orange powder, soluble in alkali with some fluorescence, dyes in red shades and does not m. 250. The K2 salt was prepared. The di-Bz-derivative m. 156 8°. Resorcinol-p-cresolphthalein, red, soluble in NaOH with green-red fluorescence less bright

than that of fluorescein, softens 220° and is precipitated from alkaline solution by CO2. It forms a K2 salt. Resorcinolfluorescein, unlike

fluorescein, insol. in Na2CO3, is soluble in NaOH with less bright fluorescence, than fluorescein, decomp. 230°.

Resorcinolisatinein, orange, is soluble in alkali with slight fluorescence,

more marked in concentrated H2SO4, dyes orange shades, and does not m. 265°. Yield 75%. Tetra-Br-derivative dyes red shades, decomp. 230°.

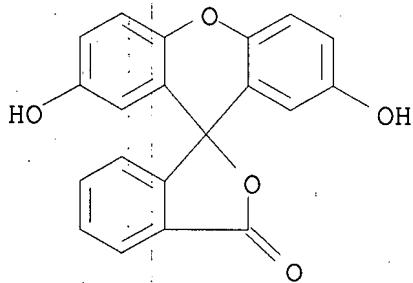
Isatinrhodamine, soluble in acids with green-red fluorescence, dyes with an impure violet shade and softens 242°. Pyrogallolisatinein gives non-fluorescent, deeply colored solns. in alkali and organic solvents, does not m. 250°. Phenolisatinein, grayish powder, soluble in alkali with red color, softens 285°. p-Cresolisatinein anhydride, soluble in glacial AcOH and concentrated H2SO4 with slight fluorescence,

does not m. 250°.

IT 596-04-3, Fluoran, 2,7-dihydroxy-
(constitution of)

RN 596-04-3 CAPLUS

CN Spiro[isobenzofuran-1(3H),9'-[9H]xanthen]-3-one, 2',7'-dihydroxy- (9CI)
(CA INDEX NAME)



=> log y

10/533,377

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
FULL ESTIMATED COST	ENTRY	SESSION
	294.83	467.14
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
CA SUBSCRIBER PRICE	ENTRY	SESSION
	-38.22	-38.22

STN INTERNATIONAL LOGOFF AT 21:32:46 ON 10 NOV 2007